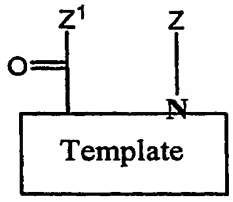
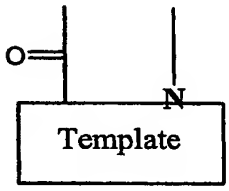
## CLAIMS

1. Compounds of the general formula

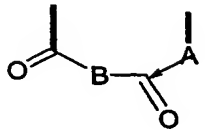



(I)

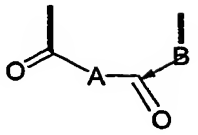
wherein



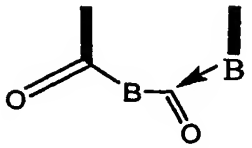

is a group of one of the formulae



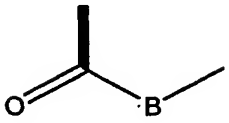

(a1)



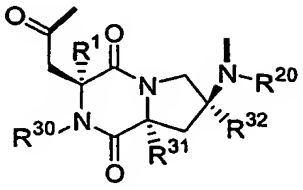

(a2)




(a3)




(a4)



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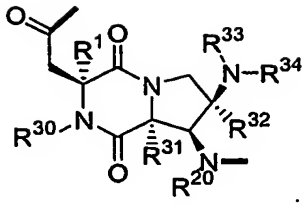
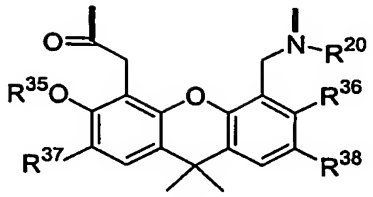
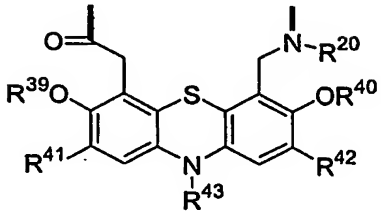


(b2)



c1)



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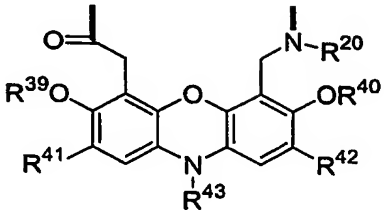
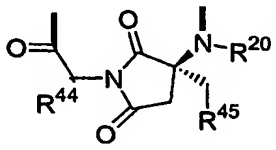


(c3)



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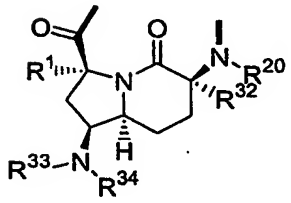
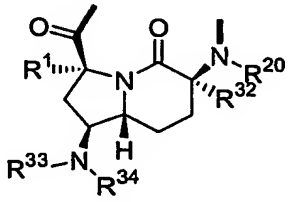


(e1)



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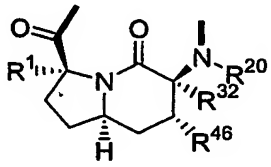

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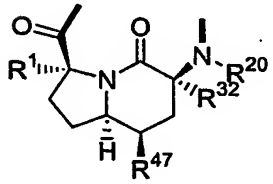

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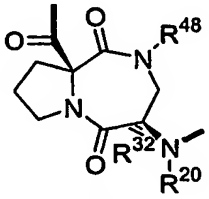
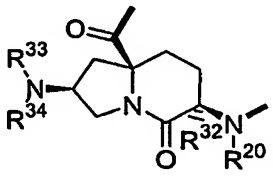
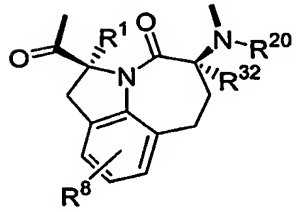


(f)



(g)



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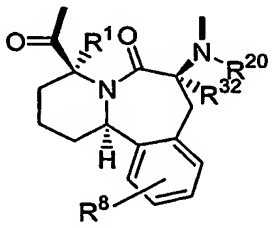
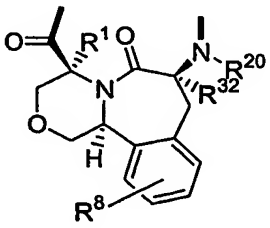


(i1)



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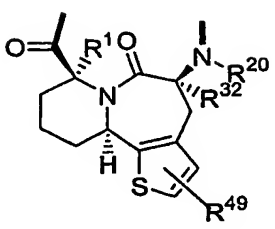

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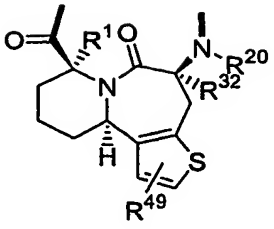

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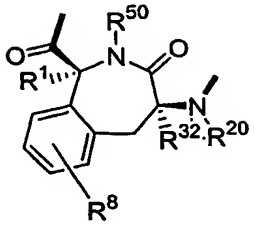

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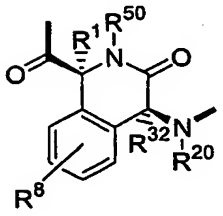

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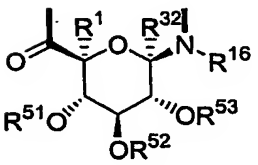

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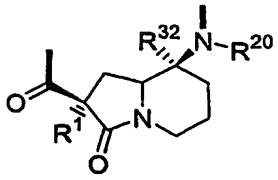

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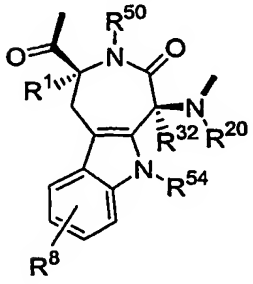
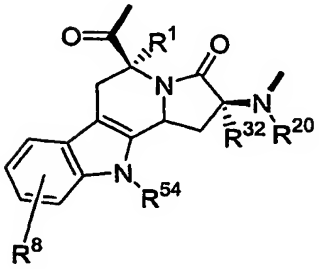
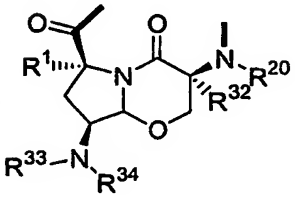


(n)



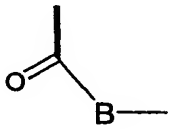
(o)

and



(p)

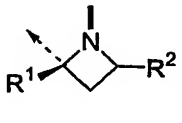
wherein



is the residue of an L-α-amino acid with B being a residue of formula $-NR^{20}CH(R^{71})-$; or the enantiomer of one of the groups **A1** to **A69** as defined hereinafter; or, in case the template is of type **(a4)**, also a residue of an amino acid with B being a residue of formula $-NR^{20}-CH_2-C_6H_4-CH_2-$;

is a group of one of the formulae

A1



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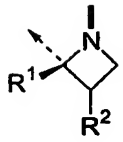

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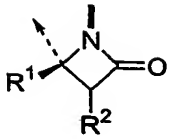

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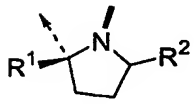

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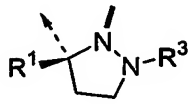

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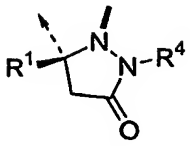

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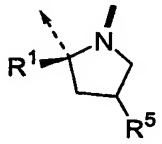

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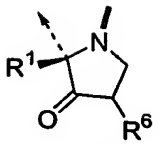

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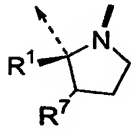

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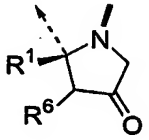

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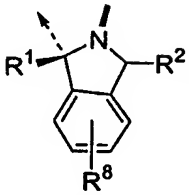

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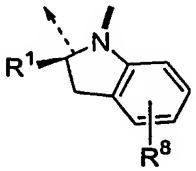

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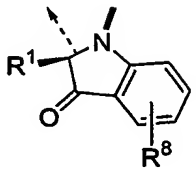

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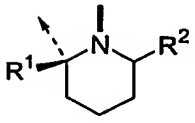
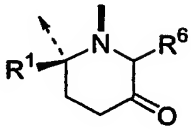


A15



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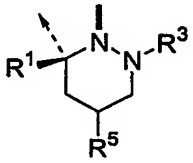

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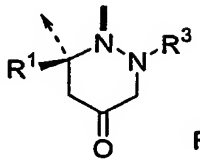

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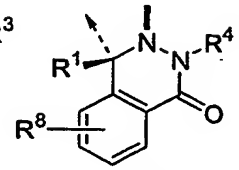

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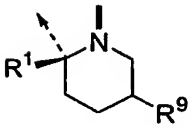

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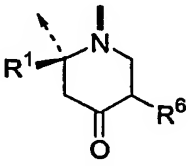

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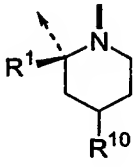

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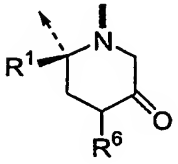

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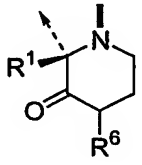

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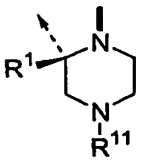

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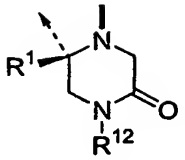

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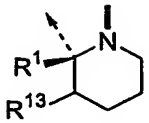

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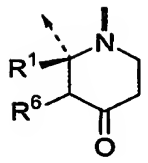

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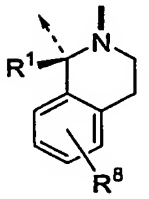

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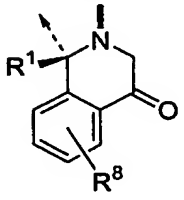

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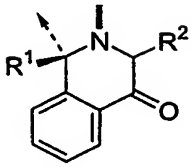

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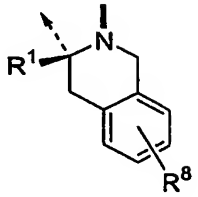

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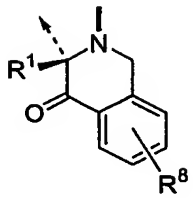

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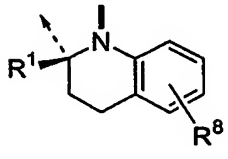

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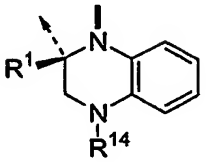

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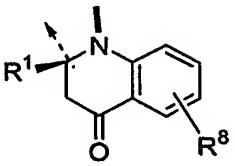

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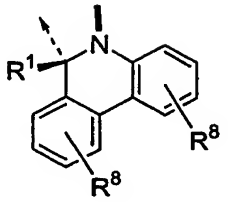

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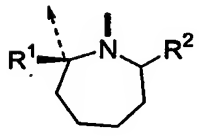

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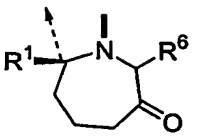

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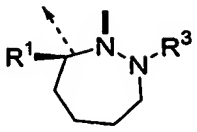

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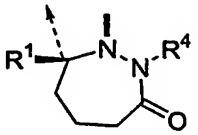

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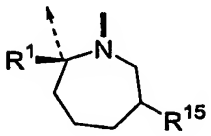

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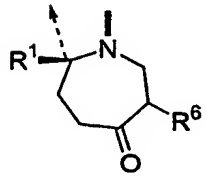

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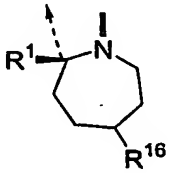

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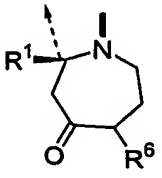

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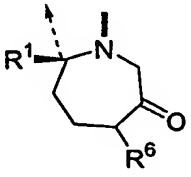

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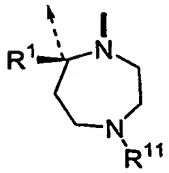

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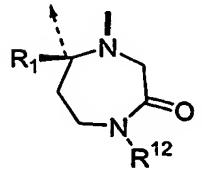

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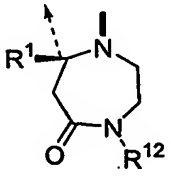

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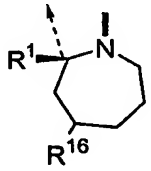

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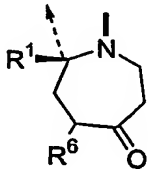

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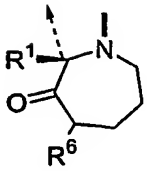

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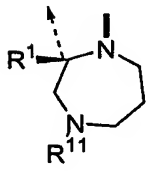

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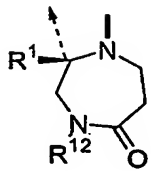

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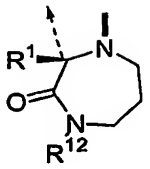

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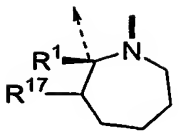

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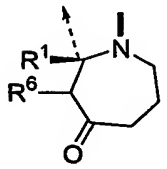

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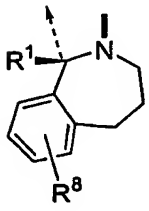

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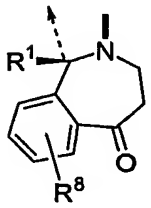

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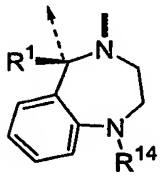

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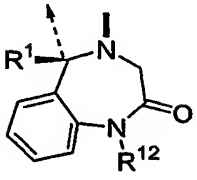

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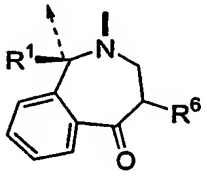

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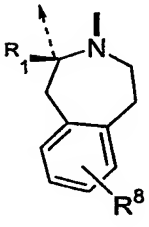

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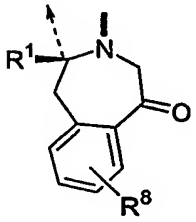

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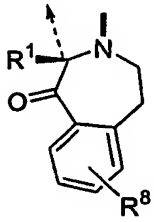

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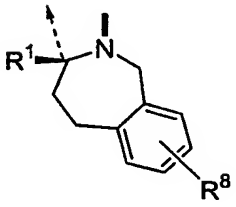

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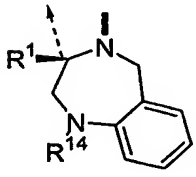

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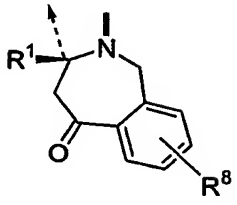

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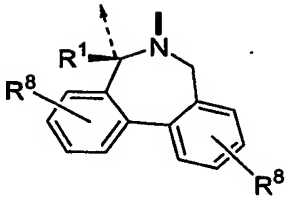

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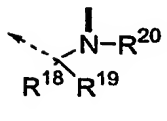

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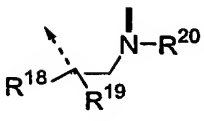

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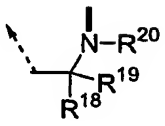

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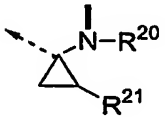

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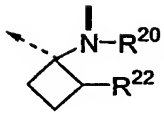

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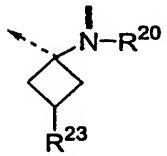

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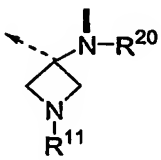

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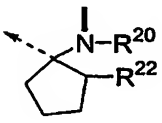

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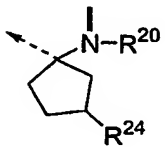

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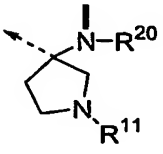

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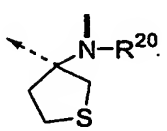

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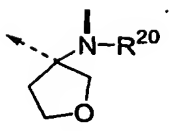

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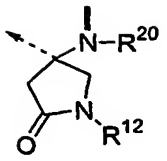

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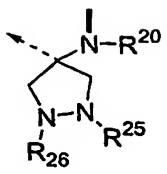

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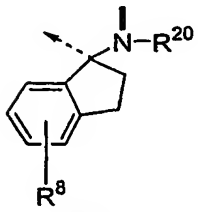

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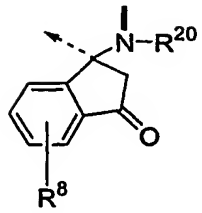

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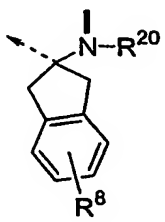

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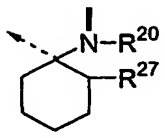

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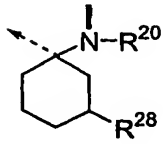

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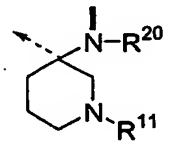

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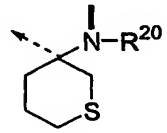

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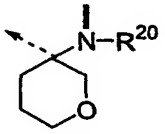

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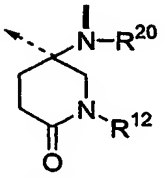

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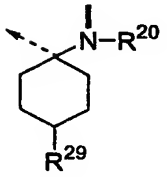

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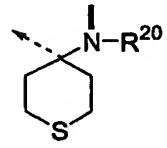
A94



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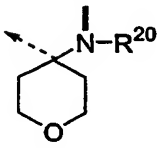

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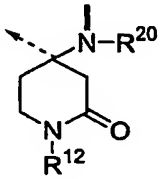

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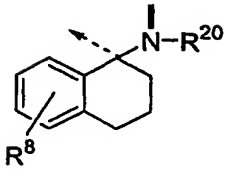

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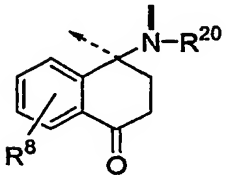

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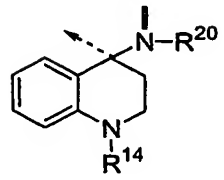

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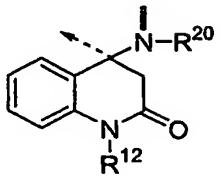

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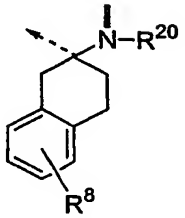

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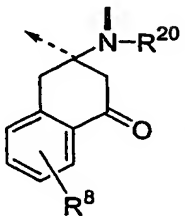


A103

and

A104

$R^1$ is H; lower alkyl; or aryl-lower alkyl;

$R^2$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$;

$-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^3$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$;
$-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$;
$-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$ ;
$-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$ ; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^4$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$; -
$(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_p(CHR^{61})_sCOOR^{57}$; $-(CH_2)_p(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_p(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_p(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^5$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^6$ is H; alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; -
$(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^7$ is alkyl; alkenyl; $-(CH_2)_q(CHR^{61})_sOR^{55}$; $-(CH_2)_q(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_q(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_q(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_r(CHR^{61})_sCOOR^{57}$; $-(CH_2)_r(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_r(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_r(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_r(CHR^{61})_s C_6H_4R^8$;

$R^8$ is H; Cl; F; $CF_3$; $NO_2$; lower alkyl; lower alkenyl; aryl; aryl-lower alkyl;
$-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})NR^{33}R^{34}$ ;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sCOR^{64}$;

$R^9$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{10}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{11}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s C_6H_4R^8$;

$R^{12}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_r(CHR^{61})_sCOOR^{57}$; $-(CH_2)_r(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_r(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_r(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_r(CHR^{61})_sC_6H_4R^8$;

$R^{13}$ is alkyl; alkenyl; $-(CH_2)_q(CHR^{61})_sOR^{55}$; $-(CH_2)_q(CHR^{61})_sSR^{56}$; $-(CH_2)_q(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_q(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_q(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_q(CHR^{61})_sCOOR^{57}$; $-(CH_2)_q(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_q(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_q(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_q(CHR^{61})_sC_6H_4R^8$;

$R^{14}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_q(CHR^{61})_sCOOR^{57}$; $-(CH_2)_q(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_q(CHR^{61})_sPO(OR^{60})_2$;. $-(CH_2)_q(CHR^{61})_sSOR^{62}$; or $-(CH_2)_q(CHR^{61})_s C_6H_4R^8$;

$R^{15}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{16}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{17}$ is alkyl; alkenyl; $-(CH_2)_q(CHR^{61})_sOR^{55}$; $-(CH_2)_q(CHR^{61})_sSR^{56}$; $-(CH_2)_q(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_q(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_q(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_q(CHR^{61})_sCOOR^{57}$; $-(CH_2)_q(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_q(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_q(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_q(CHR^{61})_sC_6H_4R^8$;

$R^{18}$ is alkyl; alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; $-(CH_2)_p(CHR^{61})_sSR^{56}$; $-(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;

$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_p(CHR^{61})_sCOOR^{57}$; $-(CH_2)_p(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_p(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_p(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{19}$ is lower alkyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; $-(CH_2)_p(CHR^{61})_sSR^{56}$; $-(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_p(CHR^{61})_sCOOR^{57}$; $-(CH_2)_p(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_p(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_p(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$; or

$R^{18}$ and $R^{19}$ taken together can form: $-(CH_2)_{2\text{-}6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or
$-(CH_2)_2NR^{57}(CH_2)_2$-;

$R^{20}$ is H; alkyl; alkenyl; or aryl-lower alkyl;

$R^{21}$ is H; alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; -
$(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{22}$ is H; alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; -
$(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{23}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{24}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{25}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$;
$-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$;
$-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$;
$-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{26}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$;

$-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$;
$-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; -
$(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$; or

$R^{25}$ and $R^{26}$ taken together can form: $-(CH_2)_{2-6}$-; $-(CH_2)_rO(CH_2)_r$-; $-(CH_2)_rS(CH_2)_r$-; or
$-(CH_2)_rNR^{57}(CH_2)_r$-;

$R^{27}$ is H; alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; -
$(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$;
$-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{28}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_s$-$OR^{55}$; $-(CH_2)_o(CHR^{61})_s SR^{56}$; $-(CH_2)_o(CHR^{61})_s$
$NR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_s COOR^{57}$; $-(CH_2)_o(CHR^{61})_s CONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_s PO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s C_6H_4R^8$;

$R^{29}$ is alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{30}$ is H; alkyl; alkenyl; or aryl-lower alkyl;

$R^{31}$ is H; alkyl; alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; $-(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s C_6H_4R^8$;

$R^{32}$ is H; lower alkyl; or aryl-lower alkyl;

$R^{33}$ is H; alkyl, alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sNR^{34}R^{63}$;
$-(CH_2)_m(CHR^{61})_sOCONR^{75}R^{82}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{78}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOR^{64}$; $-(CH_2)_o(CHR^{61})_s$-$CONR^{58}R^{59}$, $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_s SO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{34}$ is H; lower alkyl; aryl, or aryl-lower alkyl;

$R^{33}$ and $R^{34}$ taken together can form: $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or
$-(CH_2)_2NR^{57}(CH_2)_2$-;

$R^{35}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$;

$-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_p(CHR^{61})_sCOOR^{57}$; $-(CH_2)_p(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_p(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_p(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_p(CHR^{61})_s\ C_6H_4R^8$;

$R^{36}$ is H, alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_p(CHR^{61})_sCOOR^{57}$; $-(CH_2)_p(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_p(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_p(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s\ C_6H_4R^8$;

$R^{37}$ is H; F; Br; Cl; $NO_2$; $CF_3$; lower alkyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; $-(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s\ C_6H_4R^8$;

$R^{38}$ is H; F; Br; Cl; $NO_2$; $CF_3$; alkyl; alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; -
$(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_sC_6H_4R^8$;

$R^{39}$ is H; alkyl; alkenyl; or aryl-lower alkyl;

$R^{40}$ is H; alkyl; alkenyl; or aryl-lower alkyl;

$R^{41}$ is H; F; Br; Cl; $NO_2$; $CF_3$; alkyl; alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; -
$(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s\ C_6H_4R^8$;

$R^{42}$ is H; F; Br; Cl; $NO_2$; $CF_3$; alkyl; alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{55}$; -
$(CH_2)_p(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s\ C_6H_4R^8$;

$R^{43}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_sPO(OR^{60})_2$;
$-(CH_2)_o(CHR^{61})_sSO_2R^{62}$; or $-(CH_2)_o(CHR^{61})_s\ C_6H_4R^8$;

$R^{44}$ is alkyl; alkenyl; $-(CH_2)_r(CHR^{61})_sOR^{55}$; $-(CH_2)_r(CHR^{61})_sSR^{56}$; $-(CH_2)_r(CHR^{61})_sNR^{33}R^{34}$;
$-(CH_2)_r(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_r(CHR^{61})_sNR^{20}CONR^{33}R^{82}$;

$-(CH_2)_r(CHR^{61})_sCOOR^{57}$; $-(CH_2)_r(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_r(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_r(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_r(CHR^{61})_sC_6H_4R^8$;

$R^{45}$ is H; alkyl; alkenyl; $-(CH_2)_o(CHR^{61})_sOR^{55}$; $-(CH_2)_o(CHR^{61})_sSR^{56}$; $-(CH_2)_o(CHR^{61})_sNR^{33}R^{34}$;

$-(CH_2)_o(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_o(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_s(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_s(CHR^{61})_sPO(OR^{60})_2$; $-(CH_2)_s(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_s(CHR^{61})_sC_6H_4R^8$;

$R^{46}$ is H; alkyl; alkenyl; or $-(CH_2)_o(CHR^{61})_pC_6H_4R^8$;

$R^{47}$ is H; alkyl; alkenyl; or $-(CH_2)_o(CHR^{61})_sOR^{55}$;

$R^{48}$ is H; lower alkyl; lower alkenyl; or aryl-lower alkyl;

$R^{49}$ is H; alkyl; alkenyl; $-(CHR^{61})_sCOOR^{57}$; $(CHR^{61})_sCONR^{58}R^{59}$; $(CHR^{61})_sPO(OR^{60})_2$; $-(CHR^{61})_sSOR^{62}$; or $-(CHR^{61})_sC_6H_4R^8$;

$R^{50}$ is H; lower alkyl; or aryl-lower alkyl;

$R^{51}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_pPO(OR^{60})_2$; $-(CH_2)_p(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_p(CHR^{61})_sC_6H_4R^8$;

$R^{52}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_pPO(OR^{60})_2$; $-(CH_2)_p(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_p(CHR^{61})_sC_6H_4R^8$;

$R^{53}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sSR^{56}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})_sCOOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; $-(CH_2)_o(CHR^{61})_pPO(OR^{60})_2$; $-(CH_2)_p(CHR^{61})_s\ SO_2R^{62}$; or $-(CH_2)_p(CHR^{61})_sC_6H_4R^8$;

$R^{54}$ is H; alkyl; alkenyl; $-(CH_2)_m(CHR^{61})_sOR^{55}$; $-(CH_2)_m(CHR^{61})_sNR^{33}R^{34}$; $-(CH_2)_m(CHR^{61})_sOCONR^{33}R^{75}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{33}R^{82}$; $-(CH_2)_o(CHR^{61})COOR^{57}$; $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$; or $-(CH_2)_o(CHR^{61})_s\ C_6H_4R^8$;

$R^{55}$ is H; lower alkyl; lower alkenyl; aryl-lower alkyl; $-(CH_2)_m(CHR^{61})_sOR^{57}$; $-(CH_2)_m(CHR^{61})_sNR^{34}R^{63}$; $-(CH_2)_m(CHR^{61})_sOCONR^{75}R^{82}$;

$-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{78}R^{82}$; $-(CH_2)_o(CHR^{61})_s-COR^{64}$; $-(CH_2)_o(CHR^{61})COOR^{57}$; or

$-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$;

$R^{56}$ is H; lower alkyl; lower alkenyl; aryl-lower alkyl; $-(CH_2)_m(CHR^{61})_sOR^{57}$; $-(CH_2)_m(CHR^{61})_sNR^{34}R^{63}$; $-(CH_2)_m(CHR^{61})_sOCONR^{75}R^{82}$; $-(CH_2)_m(CHR^{61})_sNR^{20}CONR^{78}R^{82}$; $-(CH_2)_o(CHR^{61})_s-COR^{64}$; or $-(CH_2)_o(CHR^{61})_sCONR^{58}R^{59}$;

$R^{57}$ is H; lower alkyl; lower alkenyl; aryl lower alkyl; or heteroaryl lower alkyl;

$R^{58}$ is H; lower alkyl; lower alkenyl; aryl; heteroaryl; aryl-lower alkyl; or heteroaryl-lower alkyl;

$R^{59}$ is H; lower alkyl; lower alkenyl; aryl; heteroaryl; aryl-lower alkyl; or heteroaryl-lower alkyl; or

$R^{58}$ and $R^{59}$ taken together can form: $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$;

$R^{60}$ is H; lower alkyl; lower alkenyl; aryl; or aryl-lower alkyl;

$R^{61}$ is alkyl; alkenyl; aryl; heteroaryl; aryl-lower alkyl; heteroaryl-lower alkyl; $-(CH_2)_mOR^{55}$; $-(CH_2)_mNR^{33}R^{34}$; $-(CH_2)_mOCONR^{75}R^{82}$; $-(CH_2)_mNR^{20}CONR^{78}R^{82}$; $-(CH_2)_oCOOR^{37}$; $-(CH_2)_oNR^{58}R^{59}$; or $-(CH_2)_oPO(COR^{60})_2$;

$R^{62}$ is lower alkyl; lower alkenyl; aryl, heteroaryl; or aryl-lower alkyl;

$R^{63}$ is H; lower alkyl; lower alkenyl; aryl, heteroaryl; aryl-lower alkyl; heteroaryl-lower alkyl;

$-COR^{64}$; $-COOR^{57}$; $-CONR^{58}R^{59}$; $-SO_2R^{62}$; or $-PO(OR^{60})_2$;

$R^{34}$and $R^{63}$ taken together can form: $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$;

$R^{64}$ is H; lower alkyl; lower alkenyl; aryl; heteroaryl; aryl-lower alkyl; heteroaryl-lower alkyl;

$-(CH_2)_p(CHR^{61})_sOR^{65}$; $-(CH_2)_p(CHR^{61})_sSR^{66}$; or $-(CH_2)_p(CHR^{61})_sNR^{34}R^{63}$; $-(CH_2)_p(CHR^{61})_sOCONR^{75}R^{82}$; $-(CH_2)_p(CHR^{61})_sNR^{20}CONR^{78}R^{82}$;

$R^{65}$ is H; lower alkyl; lower alkenyl; aryl, aryl-lower alkyl; heteroaryl-lower alkyl; $-COR^{57}$; $-COOR^{57}$; or $-CONR^{58}R^{59}$;

$R^{66}$ is H; lower alkyl; lower alkenyl; aryl; aryl-lower alkyl; heteroaryl-lower alkyl; or $-CONR^{58}R^{59}$;

Z and $Z^1$ are chains of n and, respectively, n' α-amino acid residues whereby either n is 4 and n' is 6 or n is 5 and n' is 7, the positions of said amino acid residues in said chain Z being counted starting from the N-terminal amino acid and the positions of said amino acid residues in said chain $Z^1$ being counted starting from the C-terminal amino acid, whereby these amino acid residues are, depending on their position in the chains, Gly, or Pro, or of one of the types

C: $-NR^{20}CH(R^{72})CO-$;

D: $-NR^{20}CH(R^{73})CO-$;

E: $-NR^{20}CH(R^{74})CO-$;

F: $-NR^{20}CH(R^{84})CO-$; and

H: $-NR^{20}-CH(CO-)-(CH_2)_{4-7}-CH(CO-)-NR^{20}-$;

$-NR^{20}-CH(CO-)-(CH_2)_pSS(CH_2)_p-CH(CO-)-NR^{20}-$;

$-NR^{20}-CH(CO-)-(-(CH_2)_pNR^{20}CO(CH_2)_p-CH(CO-)-NR^{20}-$;

$-NR^{20}-CH(CO-)-(-(CH_2)_pNR^{20}CONR^{20}(CH_2)_p-CH(CO-)-NR^{20}-$; and

I: $-NR^{86}CH_2CO-$;

$R^{71}$ is lower alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{75}$; $-(CH_2)_p(CHR^{61})_sSR^{75}$;
$-(CH_2)_p(CHR^{61})_sOCONR^{33}R^{75}$;
$-(CH_2)_o(CHR^{61})_sCOOR^{75}$; $-(CH_2)_pCONR^{58}R^{59}$; $-(CH_2)_pPO(OR^{62})_2$; $-(CH_2)_pSO_2R^{62}$; or
$-(CH_2)_o-C_6R^{67}R^{68}R^{69}R^{70}R^{76}$;

$R^{72}$ is H, lower alkyl; lower alkenyl; $-(CH_2)_p(CHR^{61})_sOR^{85}$; or $-(CH_2)_p(CHR^{61})_sSR^{85}$;

$R^{73}$ is $-(CH_2)_oR^{77}$; $-(CH_2)_rO(CH_2)_oR^{77}$; $-(CH_2)_rS(CH_2)_oR^{77}$; or $-(CH_2)_rNR^{20}(CH_2)_oR^{77}$;

$R^{74}$ is $-(CH_2)_pNR^{78}R^{79}$; $-(CH_2)_pNR^{77}R^{80}$; $-(CH_2)_pC(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_pC(=NOR^{50})NR^{78}R^{79}$;
$-(CH_2)_pC(=NNR^{78}R^{79})NR^{78}R^{79}$; $-(CH_2)_pNR^{80}C(=NR^{80})NR^{78}R^{79}$;
$-(CH_2)_pN=C(NR^{78}R^{80})NR^{79}R^{80}$;$-(CH_2)_pC_6H_4NR^{78}R^{79}$; $-(CH_2)_pC_6H_4NR^{77}R^{80}$;
$-(CH_2)_pC_6H_4C(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_pC_6H_4C(=NOR^{50})NR^{78}R^{79}$;
$-(CH_2)_pC_6H_4C(=NNR^{78}R^{79})NR^{78}R^{79}$; $-(CH_2)_pC_6H_4NR^{80}C(=NR^{80})NR^{78}R^{79}$;
$-(CH_2)_pC_6H_4N=C(NR^{78}R^{80})NR^{79}R^{80}$; $-(CH_2)_rO(CH_2)_mNR^{78}R^{79}$; $-(CH_2)_rO(CH_2)_mNR^{77}R^{80}$;
$-(CH_2)_rO(CH_2)_pC(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_rO(CH_2)_pC(=NOR^{50})NR^{78}R^{79}$;
$-(CH_2)_rO(CH_2)_pC(=NNR^{78}R^{79})NR^{78}R^{79}$; $-(CH_2)_rO(CH_2)_mNR^{80}C(=NR^{80})NR^{78}R^{79}$;
$-(CH_2)_rO(CH_2)_mN=C(NR^{78}R^{80})NR^{79}R^{80}$; $-(CH_2)_rO(CH_2)_pC_6H_4CNR^{78}R^{79}$;
$-(CH_2)_rO(CH_2)_pC_6H_4C(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_rO(CH_2)_pC_6H_4C(=NOR^{50})NR^{78}R^{79}$;
$-(CH_2)_rO(CH_2)_pC_6H_4C(=NNR^{78}R^{79})NR^{78}R^{79}$;
$-(CH_2)_rO(CH_2)_pC_6H_4NR^{80}C(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_rS(CH_2)_mNR^{78}R^{79}$;
$-(CH_2)_rS(CH_2)_mNR^{77}R^{80}$;$-(CH_2)_rS(CH_2)_pC(=NR^{80})NR^{78}R^{79}$;

$-(CH_2)_rS(CH_2)_pC(=NOR^{50})NR^{78}R^{79}$; $-(CH_2)_rS(CH_2)_pC(=NNR^{78}R^{79})NR^{78}R^{79}$;
$-(CH_2)_rS(CH_2)_mNR^{80}C(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_rS(CH_2)_mN=C(NR^{78}R^{80})NR^{79}R^{80}$;
$-(CH_2)_rS(CH_2)_pC_6H_4CNR^{78}R^{79}$; $-(CH_2)_rS(CH_2)_pC_6H_4C(=NR^{80})NR^{78}R^{79}$;
$-(CH_2)_rS(CH_2)_pC_6H_4C(=NOR^{50})NR^{78}R^{79}$; $-(CH_2)_rS(CH_2)_pC_6H_4C(=NNR^{78}R^{79})NR^{78}R^{79}$;
$-(CH_2)_rS(CH_2)_pC_6H_4NR^{80}C(=NR^{80})NR^{78}R^{79}$; $-(CH_2)_pNR^{80}COR^{64}$; $-(CH_2)_pNR^{80}COR^{77}$;
$-(CH_2)_pNR^{80}CONR^{78}R^{79}$; or $-(CH_2)_pC_6H_4NR^{80}CONR^{78}R^{79}$;

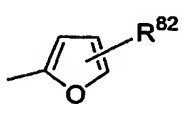
$R^{75}$ is lower alkyl; lower alkenyl; or aryl-lower alkyl;

$R^{33}$ and $R^{75}$ taken together can form: $-(CH_2)_{2\text{-}6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-;

$R^{75}$ and $R^{82}$ taken together can form: $-(CH_2)_{2\text{-}6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-;

$R^{76}$ is H; lower alkyl; lower alkenyl; aryl-lower alkyl; $-(CH_2)_oOR^{72}$; $-(CH_2)_oSR^{72}$; $-(CH_2)_oNR^{33}R^{34}$; $-(CH_2)_oOCONR^{33}R^{75}$; $-(CH_2)_oNR^{20}CONR^{33}R^{82}$; $-(CH_2)_oCOOR^{75}$; $-(CH_2)_oCONR^{58}R^{59}$; $-(CH_2)_oPO(OR^{60})_2$; $-(CH_2)_pSO_2R^{62}$; or $-(CH_2)_oCOR^{64}$;

$R^{77}$ is $-C_6R^{67}R^{68}R^{69}R^{70}R^{76}$; or a heteroaryl group of one of the formulae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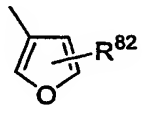
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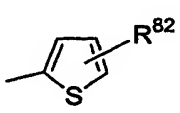
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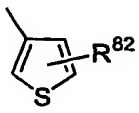
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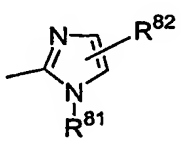
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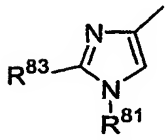
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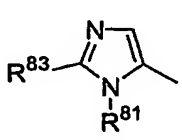
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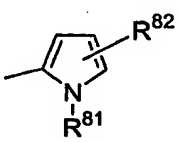
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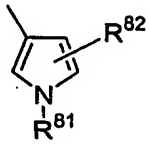
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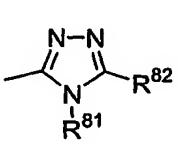
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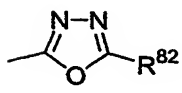
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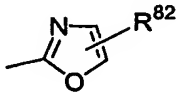
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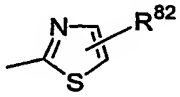
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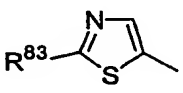
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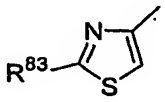
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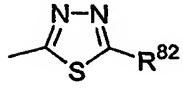
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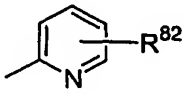
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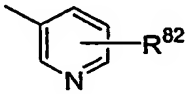
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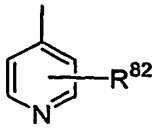
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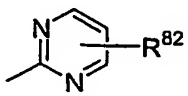
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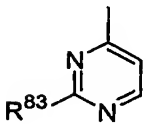

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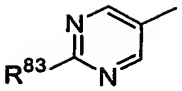

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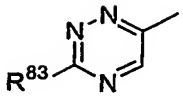

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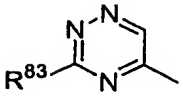

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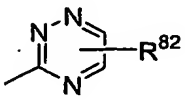

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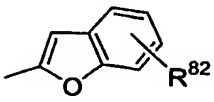

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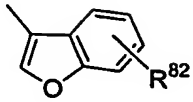

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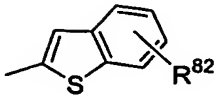

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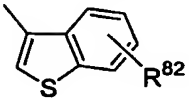

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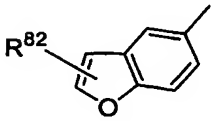

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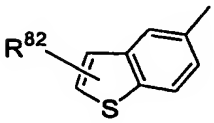

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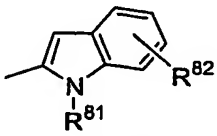

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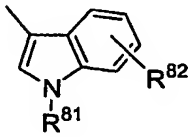

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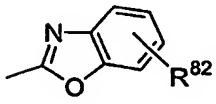

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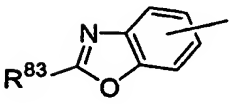

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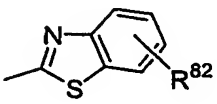

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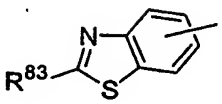

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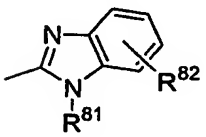

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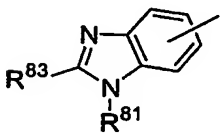

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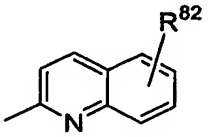

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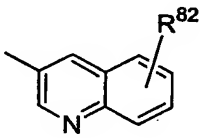

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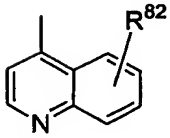

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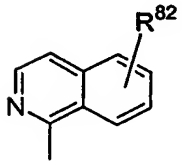

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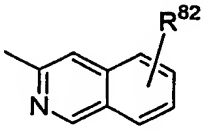

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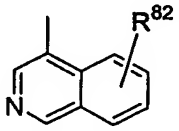

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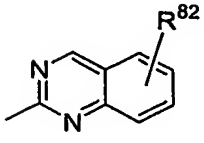

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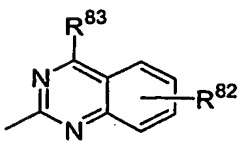

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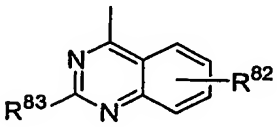

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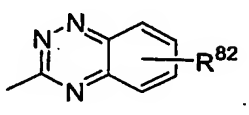


H49

$R^{78}$ is H; lower alkyl; aryl; or aryl-lower alkyl;

$R^{78}$ and $R^{82}$ taken together can form: $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$;

$R^{79}$ is H; lower alkyl; aryl; or aryl-lower alkyl; or

$R^{78}$ and $R^{79}$, taken together, can be $-(CH_2)_{2-7}-$; $-(CH_2)_2O(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$;

$R^{80}$ is H; or lower alkyl;

$R^{81}$ is H; lower alkyl; or aryl-lower alkyl;

$R^{82}$ is H; lower alkyl; aryl; heteroaryl; or aryl-lower alkyl;

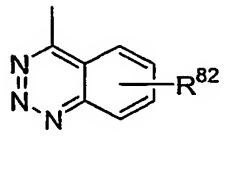
$R^{33}$ and $R^{82}$ taken together can form: $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$;

$R^{83}$ is H; lower alkyl; aryl; or $-NR^{78}R^{79}$;

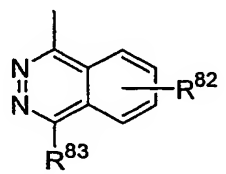
$R^{84}$ is $-(CH_2)_pCONR^{78}R^{79}$; $-(CH_2)_pNR^{80}CONR^{78}R^{79}$; $-(CH_2)_pC_6H_4CONR^{78}R^{79}$; or $-(CH_2)_pC_6H_4NR^{80}CONR^{78}R^{79}$;

$R^{85}$ is lower alkyl; or lower alkenyl;

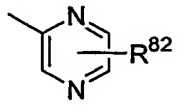
$R^{86}$ is $R^{74}$; $-[(CH_2)_u-X]_t-(CH_2)_vNR^{78}R^{79}$; $-[(CH_2)_u-X]_t-(CH_2)_v-C(=NR^{80})NR^{78}R^{79}$; X is -O-, $-NR^{20}$-,



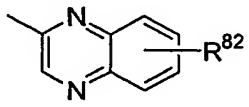
**H50**



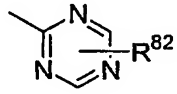
**H51**



**H52**



**H53**



**H54**

-S-, -OCOO-, u is 1-3, t is 1-6, v is 1-3;

with the proviso that in said chains Z and $Z^1$ of n and , respectively, n' α-amino acid residues

- if n is 4 and n' is 6, the amino acid residues in positions 1 to 4 of Z and in positions 1' to 6' of $Z^1$ are:

  - P1: of type C or of type D or of type E or of type F, or the residue is Pro;
  - P2: of type E or of type F;
  - P3: of type F, or the residue is Pro;
  - P4: of type E;

- P1': of type C or of type D or of type E or of type F, or the residue is Gly;
- P2': of type D or of type C;
- P3': of type F or the residue is Pro;
- P4': of type D or of type C;
- P5': of type E, or of type F or the residue is Pro; and
- P6': of type E or of type F, or the residue is Pro; or

- P3 and P3', taken together, can form a group of type H;

and

- if n is 5 and n' is 7, the amino acid residues in positions 1 to 5 of Z and in positions 1' to 7' of $Z^1$ are:

- P1: of type C or of type D or of type E or of type F, or the residue is Pro;
- P2: of type E or of type F;
- P3: of type F, or the residue is Pro;
- P4: of type F;
- P5: of type E

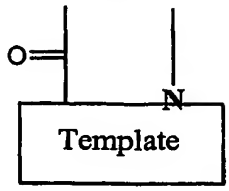
- P1': of type C or of type D or of type E or of type F, or the residue is Pro;
- P2': of type F;
- P3': of type D or the residue is Pro;
- P4': of type E or of type F;
- P5': of type D, or the residue is Pro;
- P6': of type E or of type F, or the residue is Pro; and
- P7': of type E or of type I, or the residue is Gly; or

- P2 and P2' and/or P4 and P4', taken together, can form a group of type H;

at P7' also D-isomers being possible,

and pharmaceutically acceptable salts thereof.

2. Compounds according to claim 1 wherein




is a group of formula (a1) or (a2).

3. Compounds according to claim 2 wherein A is a group of one of the formulae A1 to A69;

$R^1$ is hydrogen or lower alkyl;

$R^2$ is H; lower alkyl; lower alkenyl; $-(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-CH_2)_mSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-;

$-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

$-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

$-(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$(where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or

$-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^3$ is H; lower alkyl; lower alkenyl; $-(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_mSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

$-(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy).

$R^4$ is H; lower alkyl; lower alkenyl; $-(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_mSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mN(R^{20})COR^{64}$(where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$: isH; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy).

$R^5$ is lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower

alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;
$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are
$-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;
$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is alkyl; alkenyl; aryl; aryl-lower alkyl; or heteroaryl-lower alkyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl);
$-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy):
$R^6$ is H; lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;
$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are
$-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;
$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or

$-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^7$ is lower alkyl; lower alkenyl; $-(CH_2)_qOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_qSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_qNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2\text{-}6}-$; $-(CH_2)_2O(CH_2)_2-$;

$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_qOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

$-(CH_2)_{2\text{-}6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_qNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2\text{-}6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_qN(R^{20})COR^{64}$(where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_rCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_qCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2\text{-}6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or

$-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_rPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_rSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2\text{-}6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2\text{-}6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2\text{-}6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is

lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-;

-$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

-$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^9$ is lower alkyl; lower alkenyl; -$(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$(where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or

-$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{10}$ is lower alkyl; lower alkenyl; -$(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$: H is or lower alkyl); -$(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl;

or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$(where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{11}$ is H; lower alkyl; lower alkenyl; -$(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together ar -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mN(R^{20})COR^{64}$ (where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{12}$ is H; lower alkyl; lower alkenyl; -$(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -

$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-;

-$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

-$(CH_2)_mN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl);

-$(CH_2)_rCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_rCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-;

-$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

-$(CH_2)_rPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy):

$R^{13}$ is lower alkyl; lower alkenyl; -$(CH_2)_qOR^{55}$ (where $R^{55}$ isis lower alkyl; or lower alkenyl); -$(CH_2)_qSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_qNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_qOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_qNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_qN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_rCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_qCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_rPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_rSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{14}$ is H; lower alkyl; lower alkenyl; -$(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl is $R^{82}$: H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mN(R^{20})COR^{64}$ (where: $R^{20}$ is H; lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$;

$-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{15}$ is lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;

$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

$-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;

$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-NR^{20}COlower\ alkyl$ ($R^{20}$=H; or lower alkyl); being particularly favoured; $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl);

$-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{16}$ is lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-;

$-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

$-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-;

$-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or

$-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy); and

$R^{17}$ is lower alkyl; lower alkenyl; $-(CH_2)_qOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_qSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_qNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-;

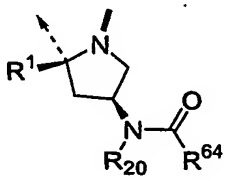
$-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_qOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

$-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_qNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}$-; $-(CH_2)_2O(CH_2)_2$-; $-(CH_2)_2S(CH_2)_2$-; or $-(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_qN(R^{20})COR^{64}$(where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_rCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_qCONR^{58}R^{59}$ (where $R^{58}$ is

lower alkyl; or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_rPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_rSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy).

4. Compounds according to claim 2 or 3 wherein A is a group of one of the formulae A5 (with $R^2$ being H); A8; A22; A25; A38 (with $R^2$ being H); A42; and A50.

5. Compounds according to claim 4 wherein A is a group of formula



**A8'**

wherein $R^{20}$ is H or lower alkyl; and $R^{64}$ is alkyl; alkenyl; aryl; aryl-lower alkyl; or heteroaryl-lower alkyl.

6. Compounds according to claim 5 wherein $R^{64}$ is n-hexyl; n-heptyl; 4-(phenyl)benzyl; diphenylmethyl, 3-amino-propyl; 5-amino-pentyl; methyl; ethyl; isopropyl; isobutyl; n-propyl; cyclohexyl; cyclohexylmethyl; n-butyl; phenyl; benzyl; (3-indolyl)methyl; 2-(3-indolyl)ethyl; (4-phenyl)phenyl; or n-nonyl.

7. Compounds according to claim 2 wherein A is a group of one of the formulae A70 to A104;
$R^{20}$ is H; or lower alkyl;
$R^{18}$ is lower alkyl;
$R^{19}$ is lower alkyl; lower alkenyl; -$(CH_2)_pOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_pSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_pNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;
-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_pOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or

$R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_pNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_pN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $(CH_2)_pCOOR^{57}$ (where $R^{57}$: lower alkyl; or lower alkenyl); $(CH_2)_pCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; or lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_pSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $(CH_2)_oC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{21}$ is H; lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl) ; $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl);

$-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{22}$ is lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower

alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;
-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are
-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$(where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl);
-$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; CF; lower alkyl; lower alkenyl; or lower alkoxy);
$R^{23}$ is H; lower alkyl; lower alkenyl; -$(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;
-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are
-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;
-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$NR^{20}$COlower alkyl ($R^{20}$=H; or lower alkyl) being particularly favoured; -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl);

$-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{24}$ is lower alkyl; lower alkenyl; $-(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;

$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

$-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$;

$-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-NR^{20}C$Olower alkyl ($R^{20}$=H ; or lower alkyl) being particularly favoured; $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl);

$-(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or $-(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{25}$ is H; lower alkyl; lower alkenyl; $-(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or

$-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or $-(CH_2)_2NR^{57}(CH_2)_2-$; where $R^{57}$ is H; or lower alkyl); $-(CH_2)_mN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); $-(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; lower alkyl; or

$R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

-$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{26}$ is H; lower alkyl; lower alkenyl; -$(CH_2)_mOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_mNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or

-$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_mN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl; or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl);

-$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy); or, alternatively, $R^{25}$ and $R^{26}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{34}(CH_2)_2$-;

$R^{27}$ is H; lower alkyl; lower alkenyl; -$(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -

$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$ (where $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2\text{-}6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy);

$R^{28}$ is lower alkyl; lower alkenyl; -$(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2\text{-}6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

-$(CH_2)_{2\text{-}6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2\text{-}6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$(where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2\text{-}6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy); and

$R^{29}$ is lower alkyl; lower alkenyl; -$(CH_2)_oOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSR^{56}$ (where $R^{56}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are -$(CH_2)_{2\text{-}6}$-; -$(CH_2)_2O(CH_2)_2$-;

-$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are

-$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oN(R^{20})COR^{64}$(where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$NR^{20}$COlower-alkyl ($R^{20}$=H; or lower alkyl) being particularly favoured; -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is lower alkyl; or lower alkenyl);
-$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl);
-$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_qC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy).

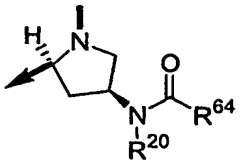
8. Compounds according to claim 7 wherein $R^{23}$, $R^{24}$ and $R^{29}$ are -$NR^{20}$-CO-lower alkyl where $R^{20}$ is H; or lower alkyl.

9. Compounds according to claim 7 or 8 wherein A is a group of one of the formulae A74 (with $R^{22}$ being H); a75; A76; A77 (with $R^{22}$ being H); A78; and A79.

10. Compounds according to any one of claims 2 to 9 wherein B is a group of formula -$NR^{20}CH(R^{71})$- or an enantiomer of one of the groups A5 (with $R^2$ being H); A8; A22; A25; A38 (with $R^2$ being H); A42; A47; and A50.

11. Compounds according to claim 10 wherein B-CO is Asn; Cys; Gln; His; Met; Phe; Pro; Ser; Thr; Trp; Tyr; Sar; 4AmPhe; 3AmPhe; 2AmPhe; Phe(mC($NH_2$)=NH; Phe(pC($NH_2$)=NH; Phe(mNHC ($NH_2$)=NH; Phe(pNHC ($NH_2$)=NH; Phg; Cha; $C_4$al; $C_5$al; 2-Nal; 1-Nal; 4Cl-Phe; 3Cl-Phe; 2Cl-Phe; 3,4$Cl_2$Phe; 4F-Phe; 3F-Phe; 2F-Phe; Tic; Thi; Tza; Mso; Y(Bzl); Bip; S(Bzl); T(Bzl); hCha; hCys; hSer; hPhe; Bpa; Pip; OctG; MePhe; MeNle; MeAla; MeIle; MeVal; MeLeu, .

12. Compounds according to claim 10 or 11 wherein B is a group, having (L)-configuration, of formula

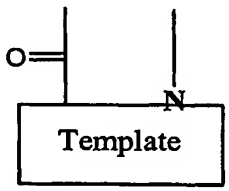


A8"

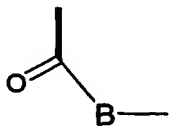
wherein $R^{20}$ is H; or lower alkyl; and $R^{64}$ is alkyl; alkenyl; aryl; aryl-lower alkyl; or heteroaryl-lower alkyl.

13. Compounds according to claim 12 wherein $R^{64}$ is n-hexyl; n-heptyl; 4-(phenyl)benzyl; diphenylmethyl, 3-amino-propyl; 5-amino-pentyl; methyl; ethyl; isopropyl; isobutyl; n-propyl; cyclohexyl; cyclohexylmethyl; n-butyl; phenyl; benzyl; (3-indolyl)methyl; 2-(3-indolyl)ethyl; (4-phenyl)phenyl; or n-nonyl.

14. Compounds according to claim 1 wherein




is a group of formula (a4) or (b1);




is the residue of AMPA;

$R^1$ is H; or lower alkyl;

$R^{20}$ is H; or lower alkyl;

$R^{30}$ is H; or methyl;

$R^{31}$ is H; lower alkyl; lower alkenyl; $-(CH_2)_pOR^{55}$ (where $R^{55}$ is lower alkyl; or lower alkenyl); $-(CH_2)_pNR^{33}R^{34}$ (where $R^{33}$ is lower alkyl; or lower alkenyl; $R^{34}$ is H; or lower alkyl; or $R^{33}$ and $R^{34}$ taken together are $-(CH_2)_{2-6}-$; $-(CH_2)_2O(CH_2)_2-$; $-(CH_2)_2S(CH_2)_2-$; or -

$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_pOCONR^{33}R^{75}$ (where $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{75}$ is lower alkyl; or $R^{33}$ and $R^{75}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_pNR^{20}CONR^{33}R^{82}$ (where $R^{20}$ is H; or lower alkyl; $R^{33}$ is H; or lower alkyl; or lower alkenyl; $R^{82}$ is H; or lower alkyl; or $R^{33}$ and $R^{82}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_pN(R^{20})COR^{64}$ (where: $R^{20}$ is H; or lower alkyl; $R^{64}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oCOOR^{57}$ (where $R^{57}$ is H; lower alkyl; or lower alkenyl); -$(CH_2)_oCONR^{58}R^{59}$ (where $R^{58}$ is lower alkyl, or lower alkenyl; and $R^{59}$ is H; lower alkyl; or $R^{58}$ and $R^{59}$ taken together are -$(CH_2)_{2-6}$-; -$(CH_2)_2O(CH_2)_2$-; -$(CH_2)_2S(CH_2)_2$-; or -$(CH_2)_2NR^{57}(CH_2)_2$-; where $R^{57}$ is H; or lower alkyl); -$(CH_2)_oPO(OR^{60})_2$ (where $R^{60}$ is lower alkyl; or lower alkenyl); -$(CH_2)_oSO_2R^{62}$ (where $R^{62}$ is lower alkyl; or lower alkenyl); or -$(CH_2)_rC_6H_4R^8$ (where $R^8$ is H; F; Cl; $CF_3$; lower alkyl; lower alkenyl; or lower alkoxy); most preferably -$CH_2CONR^{58}R^{59}$ (where $R^{58}$ is H; or lower alkyl; and $R^{59}$ is lower alkyl; or lower alkenyl); and

$R^{32}$ is H; or methyl.

15. Compounds according to claim 14 wherein $R^1$ is H; $R^{20}$ is H; $R^{30}$ is H; $R^{31}$ is carboxymethyl; or lower alkoxycarbonylmethyl; and $R^{32}$ is H.

16. Compounds according to any one of claims 1 to 15 wherein n is 4, n' is 6 and the α-amino acid residues in positions 1 to 4 of the chain Z and 1'-6' in chain $Z^1$ are:

- P1: of type D or of type E or of type F, or the residue is Pro;
- P2: of type E or of type F;
- P3: of type F, or the residue is Pro;
- P4: of type E;

- P1': of type E or of type F, or the residue is Gly;
- P2': of type D;
- P3': of type F or the residue is Pro;
- P4': of type D;
- P5': of type E, or of type F or the residue is Pro; and
- P6': of type E or of type F, or the residue is Pro; or
- P3 and P3', taken together, can form a group of type H

17. Compounds according to any one of claims 1 to 15 wherein n is 5, n' is 7 and the α-amino acid residues in positions 1 to 5 of the chain Z and 1'-7' in chain $Z^1$ are:

- P1: of type D or of type E or of type F, or the residue is Pro;
- P2: of type E or of type F;
- P3: of type F, or the residue is Pro;
- P4: of type F;
- P5: of type E

- P1': of type D or of type E or of type F, or the residue is Pro;
- P2': of type F;
- P3': of type D or the residue is Pro;
- P4': of type F;
- P5': of type D, or the residue is Pro;
- P6': of type E or of type F, or the residue is Pro; and
- P7': of type E or of type I, or the residue is Gly; or
- P2 and P2' and/or P4 and P4', taken together, can form a group of type H;

at P7`also D-isomers being possible.

18. Compounds according to claim 16 wherein the α-amino acid residues in positions 1 to 4 of the chain Z and the α-amino acid residues in positions 1' to 6' chain $Z^1$ are:

- P1: Tyr, or Arg;
- P2: Cit, or Arg;
- P3: Cys;
- P4: Arg-$NH_2$;
- P1': Lys, or Arg;
- P2': Tyr;
- P3': Cys;
- P4': 2-Nal;
- P5': Arg; and
- P6': Arg.
- Cys at pos P3 and P3` form a disulfide bridge

19. Compounds according to claim 17 wherein the α-amino acid residues in positions 1 to 5 of the chain Z and the α-amino acid residues in positions 1' to 7' chain $Z^I$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg, or Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal, Trp, F(p$NH_2$), or W(6-Cl);
- P6': Arg; and
- P7`: $^D$Arg, Arg, Ac-Arg, iPr-Arg, (EA)G, (PrA)G, (BA)G, (EGU)G,
- (PrGU)G, or (BGU)G.
- Cys at pos P4 and P4` form a disulfide bridge

20. A compound of formula I according to claim 1 wherein the template is $^L$Lys-$^L$Pro, n is 4, n' is 6 and the amino acid residues in positions 1 to 4 of the chain Z and the amino acid residues in positions 1' to 6' chain $Z^I$ are :

- P1: Tyr;
- P2: Cit;
- P3: Cys;
- P4: Arg-$NH_2$;
- P1': Arg;
- P2': Tyr;
- P3': Cys;
- P4': 2-Nal;
- P5': Arg; and
- P6': Arg.

Cys at position P4` and P4 are linked by a disulfide bridge

21. A compound of formula I according to claim 1 wherein the template is $^{D}$Pro-$^{L}$Pro, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^1$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal;
- P6': Arg; and
- P7': Arg.

Cys at position P4` and P4 form a disulfide bridge

22. A compound of formula I according to claim 1 wherein the template is $^{D}$Pro-$^{L}$Pro, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^1$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal;
- P6': Arg; and
- P7': Ac-Arg.

Cys at position P4` and P4 form a disulfide bridge

23. A compound of formula I according to claim 1 wherein the template is $^{D}$Pro-$^{L}$Pro, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^{1}$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal
- P6': Arg; and
- P7': $^{D}$Arg.

Cys at position P4` and P4 form a disulfide bridge

24. A compound of formula I according to claim 1 wherein the template is $^{D}$Pro-$^{L}$Pro, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^{1}$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': Phe(p$NH_2$);
- P6': Arg; and
- P7': Arg.

Cys at position P4` and P4 form a disulfide bridge

25. A compound of formula I according to claim 1 wherein the template is $^{D}$Pro-$^{L}$Pro, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^{1}$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal;
- P6': Arg; and
- P7': (PrA)G.

Cys at position P4` and P4 form a disulfide bridge

26. A compound of formula I according to claim 1 wherein the template is $^{D}$Pro-$^{L}$Pro, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^{1}$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal;
- P6': Arg; and
- P7': Arg.

Cys at position P4` and P4 form a disulfide bridge

27. A compound of formula I according to claim 1 wherein the template is (b1)-154, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^1$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal;
- P6': Arg; and
- P7': Arg.

Cys at position P4` and P4 form a disulfide bridge

28. A compound of formula I according to claim 1 wherein the template is AMPA, n is 5, n' is 7 and the amino acid residues in positions 1 to 5 of the chain Z and the amino acid residues in positions 1' to 7' chain $Z^1$ are:

- P1: Tyr;
- P2: Arg;
- P3: Cit;
- P4: Cys;
- P5: Arg-$NH_2$;
- P1': Lys;
- P2': Cit;
- P3': Tyr;
- P4': Cys;
- P5': 2-Nal;
- P6': Arg; and
- P7': Arg.

Cys at position P4` and P4 form a disulfide bridge

29. Enantiomers of the compounds of formulae I as defined in claim 1.

30. Compounds according to any one of claims 1 to 29 for use as therapeutically active substances.

31. Compounds according the claims 29 for use as CXCR4 antagonists.

32. A pharmaceutical composition containing a compound according to any one of claims 1 to 29 and a pharmaceutically inert carrier.

33. Compositions according to claim 32 in a form suitable for oral, topical, transdermal, injection, buccal, transmucosal, pulmonary or inhalation administration.

34. Compositions according to claim 32 or 33 in form of tablets, dragees, capsules, solutions, liquids, gels, plaster, creams, ointments, syrup, slurries, suspensions, spray, nebuliser or suppositories.

35. The use of compounds according to any one of claims 1 to 29 for the manufacture of a medicament for treating or preventing of HIV infections, or for treatment of cancer or for treatment of inflammatory disorders.

36. A process for the manufacture of compounds according to any one of claims 1-28 which process comprises

(a) coupling an appropriately functionalized solid support with an appropriately N-protected derivative of that amino acid which in the desired end-product is in position 4 of Z if n is 4 or in position 5 of Z if n is 5, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

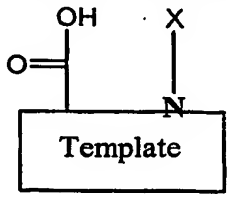
(b) removing the N-protecting group from the product thus obtained;

(c) coupling the product thus obtained with an appropriately N-protected derivative of that amino acid which in Z of the desired end-product is one position nearer the N-terminal amino acid residue, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

(d) removing the N-protecting group from the product thus obtained;

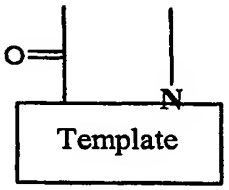
(e) repeating steps (c) and (d) until the N-terminal amino acid residue of Z has been introduced;

(f) coupling the product thus obtained with a compound of the general formula

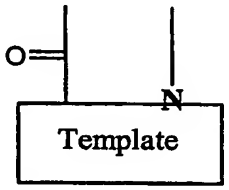



II

wherein




is as defined in claim 1 and X is an N-protecting group or, if




is to be group (a1), or (a2), above, alternatively

(fa) coupling the product obtained in step (e) with an appropriately N-protected derivative of an amino acid of the general formula

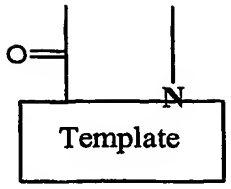
HOOC-B-H III or HOOC-A-H IV

wherein B and A are as defined in claim 1, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

(fb) removing the N-protecting group from the product thus obtained; and

(fc) coupling the product thus obtained with an appropriately N-protected derivative of an amino acid of the above general formula IV and, respectively, III, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected; or

if




is to be group (a3), above, alternatively

(fa') coupling the product obtained in step (e) with an appropriately N-protected derivative of an amino acid of the above general formula III, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

(fb') removing the N-protecting group from the product thus obtained; and

(fc') coupling the product thus obtained with an appropriately N-protected derivative of an amino acid of the above general formula III, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

(g) removing the N-protecting group from the product obtained in step (f) or (fc) or (fc');

(h) coupling the product thus obtained with an appropriately N-protected derivative of that amino acid which in the desired end-product is in position 1 of $Z^1$, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

(i) removing the N-protecting group from the product thus obtained;

(j) coupling the product thus obtained with an appropriately N-protected derivative of that amino acid which in the desired end-product is one position farther away from position 1 of $Z^1$, any functional group which may be present in said N-protected amino acid derivative being likewise appropriately protected;

(k) removing the N-protecting group from the product thus obtained;

(l) repeating steps (j) and (k) until all amino acid residues of $Z^1$ have been introduced;

(m) if desired, selectively deprotecting one or several protected functional group(s) present in the molecule and appropriately substituting the reactive group(s) thus liberated;

(n) if desired, forming one or two interstrand linkage(s) between side-chains of appropriate amino acid residues at opposite positions of the β-strand region;

(o) detaching the product thus obtained from the solid support and removing any protecting groups present on functional groups of any members of the chain of amino acid residues and, if desired, any protecting group(s) which may in addition be present in the molecule; and

(p) if desired, converting the product thus obtained into a pharmaceutically acceptable salt or converting a pharmaceutically acceptable, or unacceptable, salt thus obtained into the corresponding free compound of formula I or into a different, pharmaceutically acceptable, salt..

37. A process according to claim 36 but wherein an amino acid residue of type I is introduced by coupling with a leaving group-containing acetylating agent, followed by nucleophilic displacement with an amine of the formula $H_2NR^{86}$ which, if necessary, is appropriately protected.

38. A process according to claim 37 wherein said leaving group-containing acetylating agent is bromo, chloro or iodo acetic acid.

39. A modification of the process according to any one of claims 36 to 38 for the manufacture of compounds according to claim 29 in which enantiomers of all chiral starting materials are used.